

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111918\
 Data File : BF110864.D
 Acq On : 19 Nov 2018 20:15
 Operator : JU/SJ
 Sample : J5993-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20180929-AMENDED-WCC-2-COMP

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.228	20	24	33	rVB	86407	124509	9.19%	0.498%
2	5.075	502	508	511	rBV	59028	76863	5.67%	0.307%
3	5.175	519	525	530	rBV	773638	1034831	76.36%	4.137%
4	5.346	550	554	559	rBV2	52898	86313	6.37%	0.345%
5	5.563	588	591	600	rVV	392179	462667	34.14%	1.850%
6	5.793	626	630	637	rBV4	64656	87560	6.46%	0.350%
7	5.957	652	658	661	rBV3	56471	88926	6.56%	0.356%
8	6.175	691	695	701	rBV2	164197	304442	22.47%	1.217%
9	6.228	701	704	707	rVB2	98288	95083	7.02%	0.380%
10	6.363	723	727	731	rBV3	76541	120256	8.87%	0.481%
11	6.416	731	736	740	rVB4	96405	135104	9.97%	0.540%
12	6.469	740	745	748	rBV4	110816	183454	13.54%	0.733%
13	6.569	759	762	767	rBV	1111107	1114485	82.24%	4.456%
14	6.669	775	779	783	rBV3	66680	121928	9.00%	0.487%
15	6.710	783	786	791	rVB	1074211	919580	67.86%	3.676%
16	6.757	791	794	798	rBV3	100657	125813	9.28%	0.503%
17	6.940	821	825	829	rVB2	478424	457062	33.73%	1.827%
18	7.045	839	843	845	rBV3	91437	108096	7.98%	0.432%
19	7.069	845	847	849	rVV2	109410	112108	8.27%	0.448%
20	7.098	849	852	859	rVB	1166479	1072817	79.17%	4.289%
21	7.198	863	869	872	rBV2	225286	268085	19.78%	1.072%
22	7.257	875	879	884	rVB2	146920	185567	13.69%	0.742%
23	7.351	893	895	900	rVB2	77993	107935	7.96%	0.432%
24	7.422	904	907	910	rVV2	109145	164794	12.16%	0.659%
25	7.463	912	914	918	rVV2	152269	146900	10.84%	0.587%
26	7.510	918	922	928	rVV	913056	970419	71.61%	3.880%
27	7.557	928	930	931	rVV	75484	71488	5.28%	0.286%
28	7.575	931	933	936	rVV3	122554	111816	8.25%	0.447%
29	7.622	936	941	944	rVV3	112879	170468	12.58%	0.681%
30	7.669	946	949	952	rVV2	72115	100219	7.40%	0.401%
31	7.698	952	954	957	rVV	158755	170888	12.61%	0.683%
32	7.740	957	961	966	rVB3	122268	214370	15.82%	0.857%
33	7.857	978	981	984	rBV3	83461	108389	8.00%	0.433%
34	7.951	994	997	1000	rBV2	130464	124770	9.21%	0.499%

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Integrator: RTE

Smoothing : ON

Sampling : 1

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Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	8.028	1006	1010	1014	rVB3	103618	135005	9.96%	0.540%
36	8.081	1016	1019	1022	rVB3	139025	130173	9.61%	0.520%
37	8.163	1031	1033	1035	rBV2	82729	87677	6.47%	0.351%
38	8.222	1041	1043	1047	rBV	431004	395581	29.19%	1.581%
39	8.263	1047	1050	1053	rVB	490652	405144	29.90%	1.620%
40	8.298	1053	1056	1061	rBV3	131472	178000	13.14%	0.712%
41	8.416	1074	1076	1079	rVB2	93068	72261	5.33%	0.289%
42	8.498	1087	1090	1095	rVB4	258593	347333	25.63%	1.389%
43	8.545	1095	1098	1102	rBV5	102467	172008	12.69%	0.688%
44	8.622	1109	1111	1114	rVB	364946	264311	19.50%	1.057%
45	8.651	1114	1116	1119	rBV2	95935	89911	6.63%	0.359%
46	8.722	1125	1128	1130	rBV3	96583	85716	6.33%	0.343%
47	8.751	1130	1133	1135	rVV3	92430	92946	6.86%	0.372%
48	8.798	1139	1141	1145	rVV4	74100	81925	6.05%	0.328%
49	8.892	1154	1157	1163	rVB3	131037	167690	12.37%	0.670%
50	8.939	1163	1165	1167	rBV	73546	73314	5.41%	0.293%
51	8.992	1171	1174	1176	rBV4	74253	92397	6.82%	0.369%
52	9.039	1180	1182	1187	rBV4	73111	77365	5.71%	0.309%
53	9.122	1193	1196	1199	rVB3	120140	148368	10.95%	0.593%
54	9.157	1199	1202	1206	rBV5	91564	112476	8.30%	0.450%
55	9.204	1206	1210	1211	rBV2	121329	128556	9.49%	0.514%
56	9.228	1211	1214	1216	rVB	292590	239837	17.70%	0.959%
57	9.298	1222	1226	1229	rBV	1599990	1355123	100.00%	5.418%
58	9.363	1234	1237	1242	rBV2	194468	266803	19.69%	1.067%
59	9.569	1267	1272	1276	rBV3	130103	216843	16.00%	0.867%
60	9.639	1281	1284	1287	rVB	187286	164345	12.13%	0.657%
61	9.681	1287	1291	1294	rBV3	723622	698225	51.52%	2.791%
62	9.775	1303	1307	1312	rVB6	115345	217686	16.06%	0.870%
63	9.839	1315	1318	1320	rBV2	189092	151459	11.18%	0.606%
64	9.886	1323	1326	1329	rVB2	196227	186316	13.75%	0.745%
65	9.928	1331	1333	1336	rBV2	86805	103916	7.67%	0.415%
66	9.957	1336	1338	1340	rBV2	130866	106919	7.89%	0.427%
67	9.986	1340	1343	1347	rBV	378900	364416	26.89%	1.457%
68	10.086	1357	1360	1363	rVB2	85303	103822	7.66%	0.415%
69	10.128	1363	1367	1369	rBV4	79425	102506	7.56%	0.410%
70	10.181	1373	1376	1379	rBV2	173565	164575	12.14%	0.658%
71	10.222	1379	1383	1386	rVV3	177104	220770	16.29%	0.883%

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72	10.298	1393	1396	1399	rBV	177826	226064	16.68%	0.904%
73	10.328	1399	1401	1406	rVB3	171076	182033	13.43%	0.728%
74	10.386	1407	1411	1413	rBV4	254948	277628	20.49%	1.110%
75	10.551	1437	1439	1443	rVB3	106696	101111	7.46%	0.404%
76	10.610	1445	1449	1452	rBV	416829	482105	35.58%	1.927%
77	10.733	1468	1470	1474	rVB4	112216	121709	8.98%	0.487%
78	10.781	1474	1478	1480	rBV4	218994	264779	19.54%	1.059%
79	10.881	1490	1495	1503	rBV2	645100	896899	66.19%	3.586%
80	11.116	1533	1535	1538	rBV3	123921	146138	10.78%	0.584%
81	11.345	1571	1574	1576	rBV	413196	338687	24.99%	1.354%
82	11.480	1594	1597	1599	rBV	350882	316214	23.33%	1.264%
83	11.504	1599	1601	1609	rVB	170074	230015	16.97%	0.920%
84	11.986	1680	1683	1686	rBV2	142835	199461	14.72%	0.797%
85	12.092	1699	1701	1704	rVV3	132072	125574	9.27%	0.502%
86	12.122	1704	1706	1709	rVB2	122346	95186	7.02%	0.381%
87	12.463	1761	1764	1766	rBV	140866	125034	9.23%	0.500%
88	12.486	1766	1768	1774	rVB4	120372	135328	9.99%	0.541%
89	12.539	1774	1777	1779	rBV	281246	274448	20.25%	1.097%
90	12.622	1788	1791	1793	rBV	93897	96038	7.09%	0.384%
91	12.704	1802	1805	1807	rBV	105071	114161	8.42%	0.456%
92	12.874	1832	1834	1838	rVB2	108384	101749	7.51%	0.407%
93	12.939	1842	1845	1849	rVB2	226517	254502	18.78%	1.017%
94	12.980	1849	1852	1856	rVB	155296	156977	11.58%	0.628%
95	13.063	1863	1866	1869	rBV	1197487	1009518	74.50%	4.036%
96	13.463	1932	1934	1936	rBV	88080	78623	5.80%	0.314%
97	13.939	2012	2015	2021	rVB2	147880	197493	14.57%	0.790%
98	14.133	2045	2048	2052	rBV	311044	346413	25.56%	1.385%
99	14.557	2118	2120	2122	rBV	85167	72189	5.33%	0.289%
100	15.668	2306	2309	2320	rVB	181047	297871	21.98%	1.191%

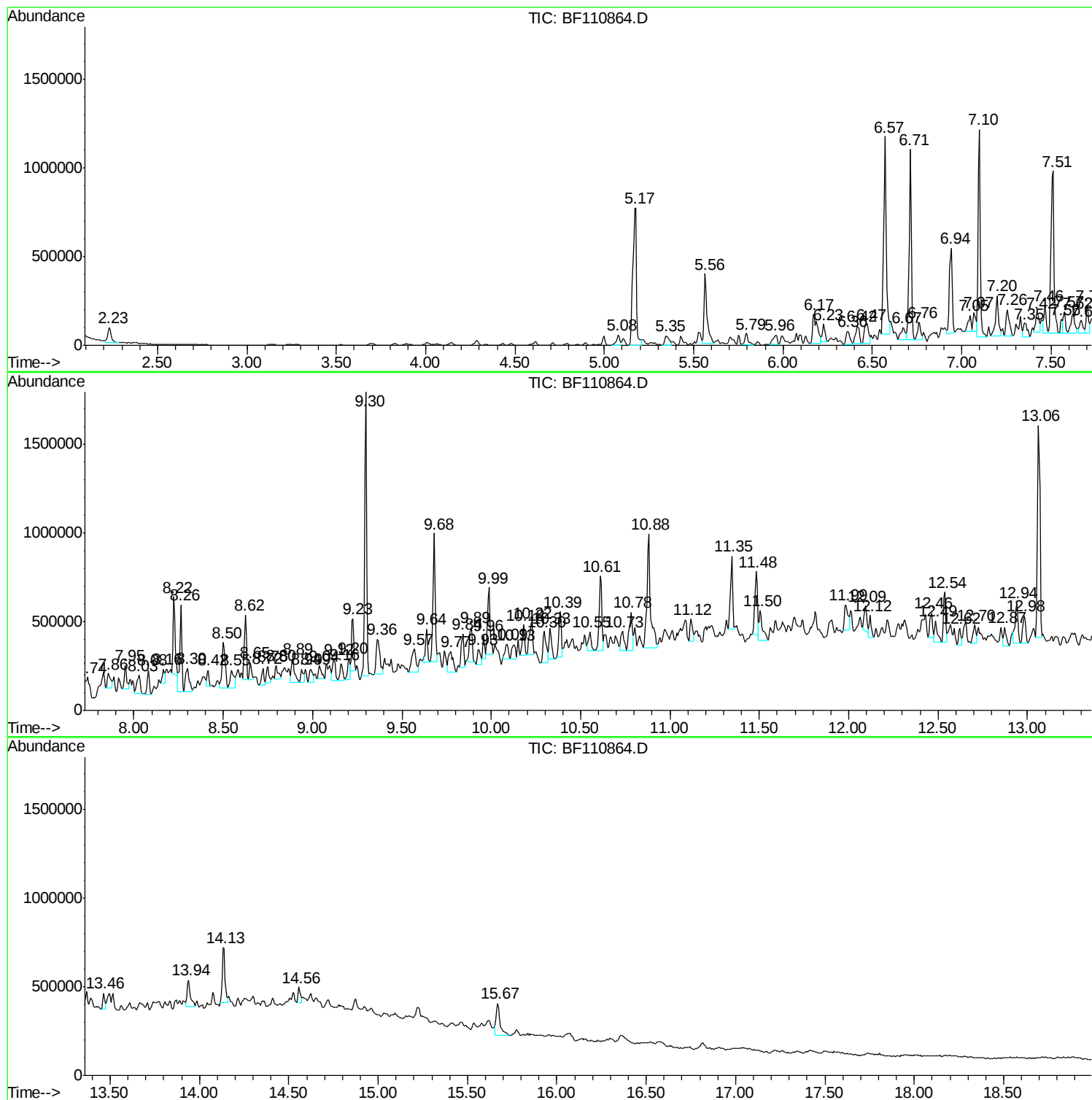
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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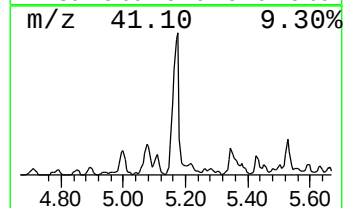
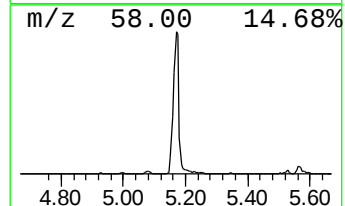
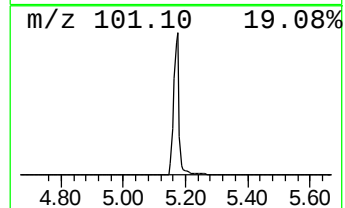
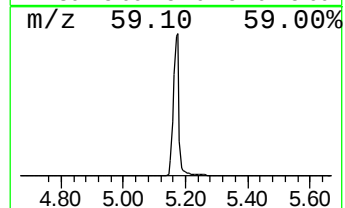
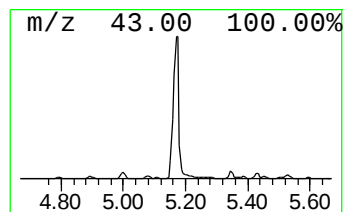
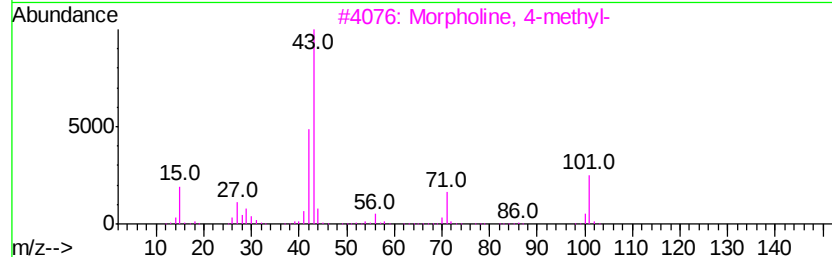
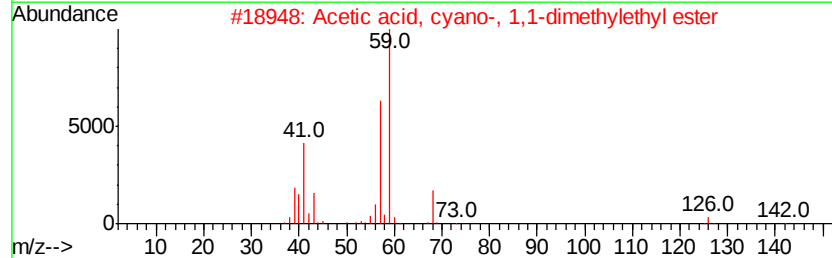
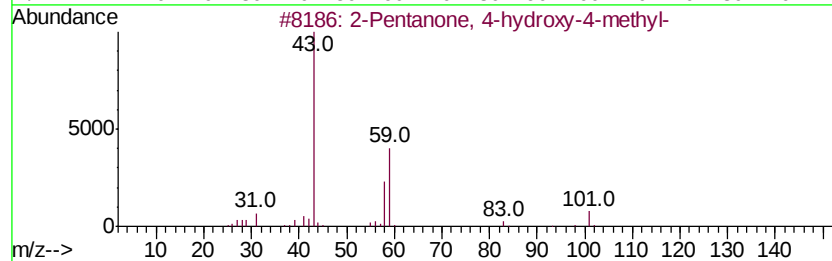
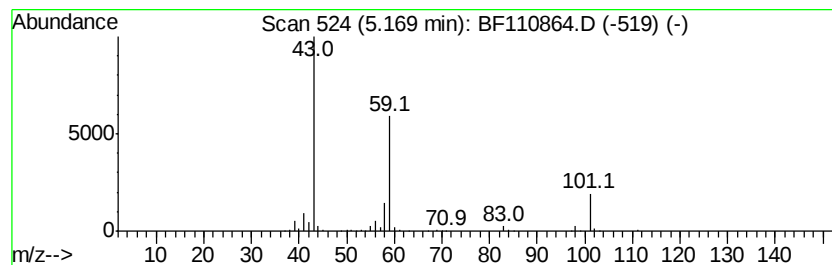
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	45.28 ng	1034830	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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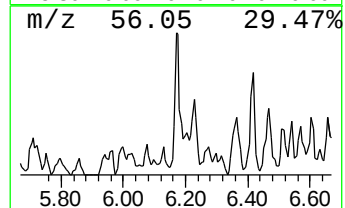
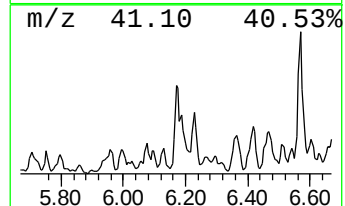
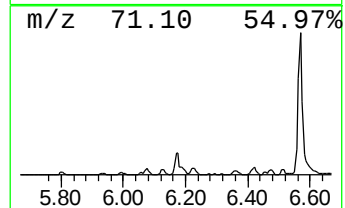
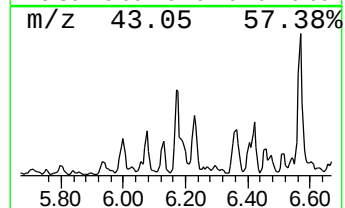
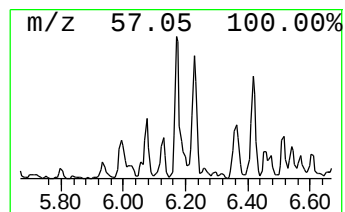
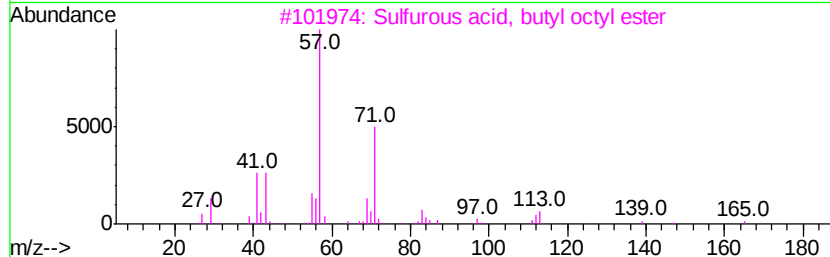
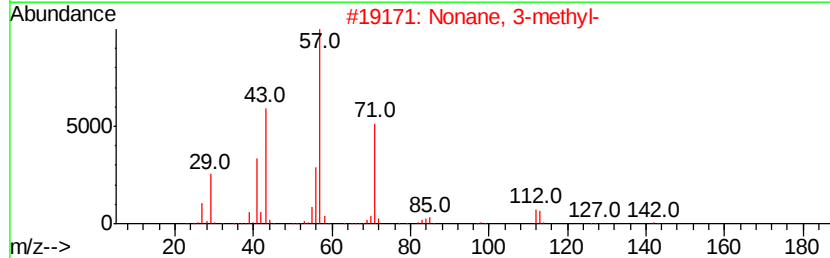
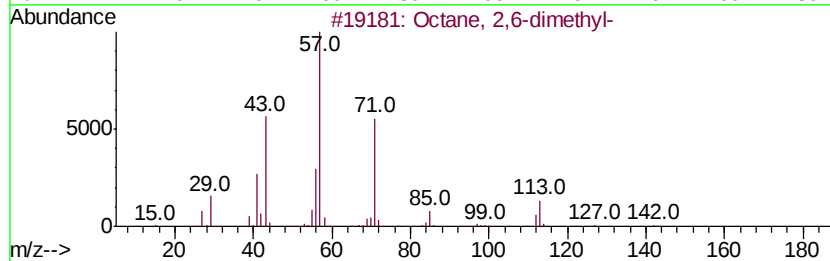
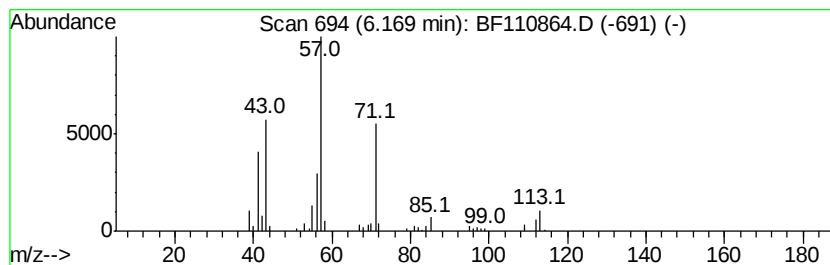
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Octane, 2,6-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.17	13.32 ng	304442	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
2		Nonane, 3-methyl-	142	C10H22	005911-04-6	64
3		Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	64
4		Sulfurous acid, butyl 2-ethylhex...	250	C12H26O3S	1000309-18-8	53
5		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	53



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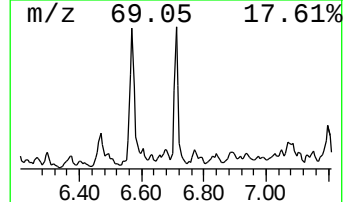
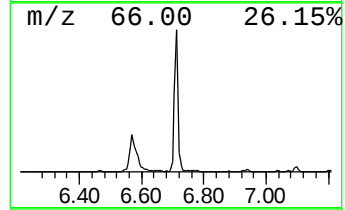
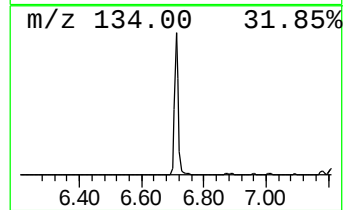
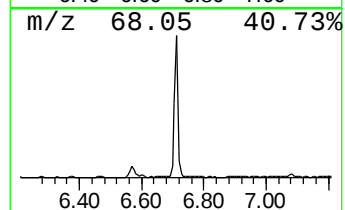
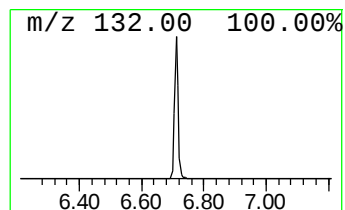
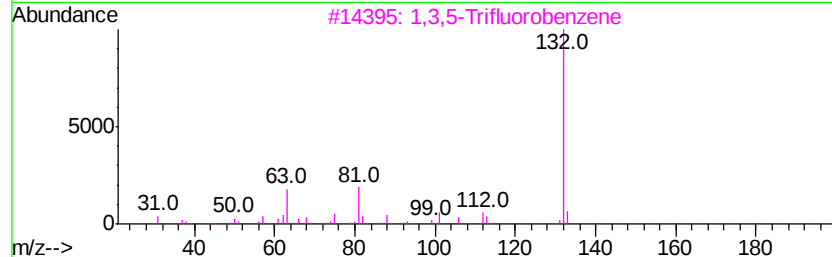
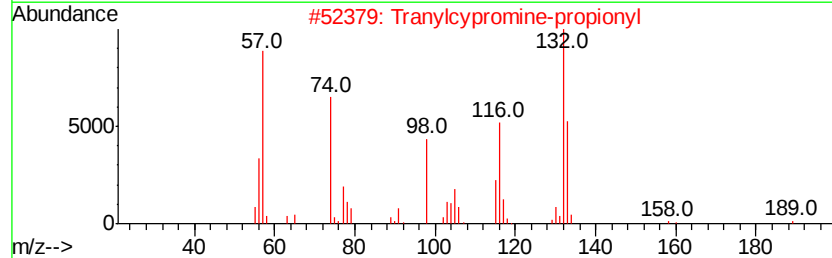
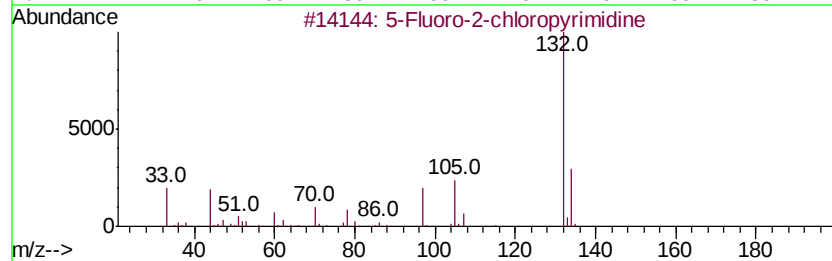
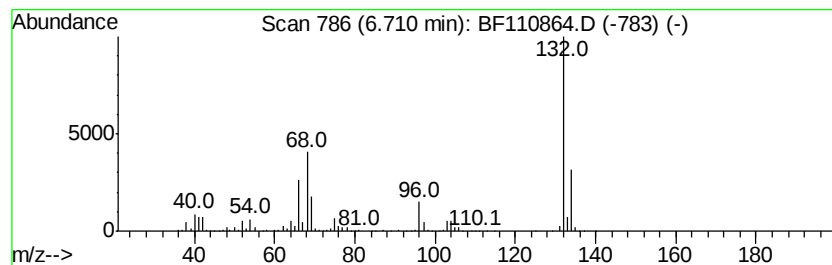
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.71 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	40.24 ng	919580	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110864.D
Acq On : 19 Nov 2018 20:15
Operator : JU/SJ
Sample : J5993-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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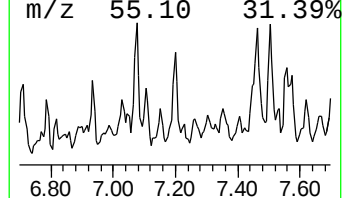
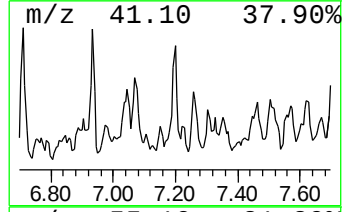
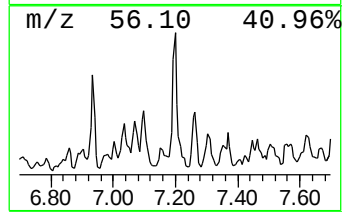
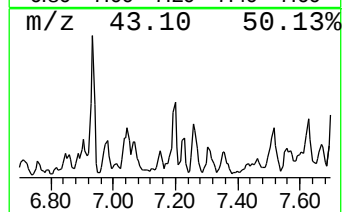
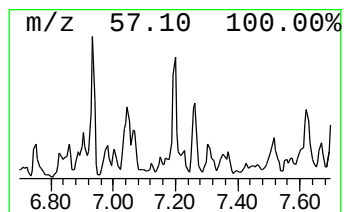
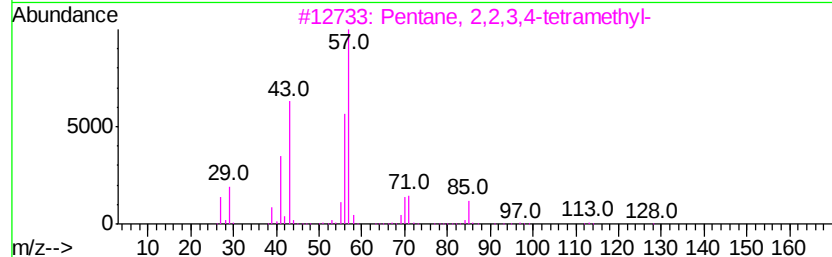
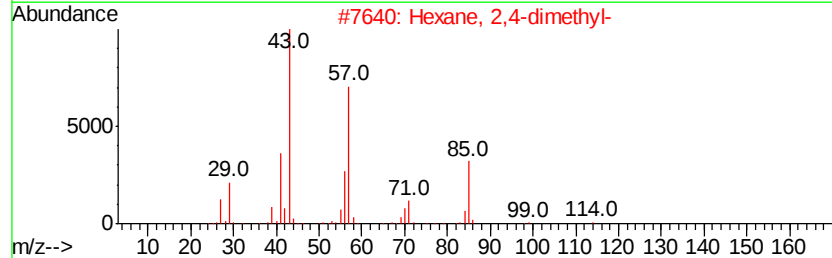
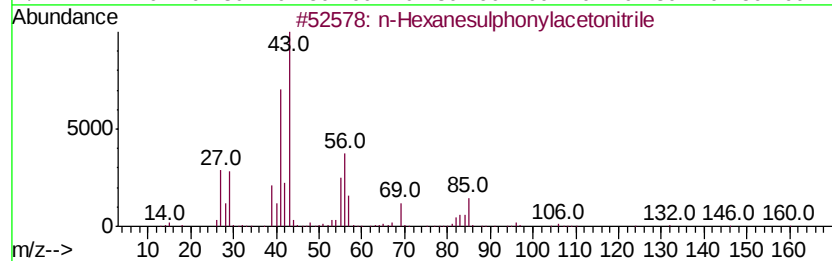
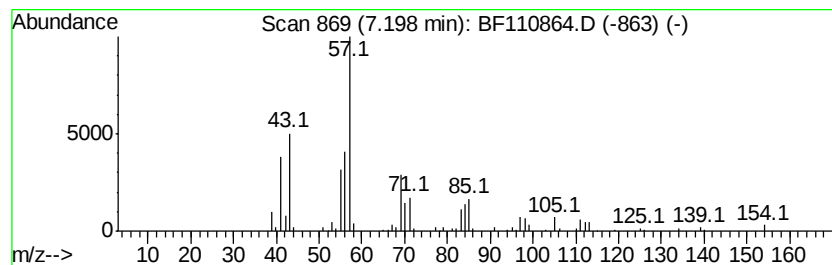
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexanesulphonylacetonitrile Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	11.73 ng	268085	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexanesulphonylacetonitrile	189	C8H15NO2S	203310-42-3	53
2		Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	50
3		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	47
4		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	47
5		Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	47



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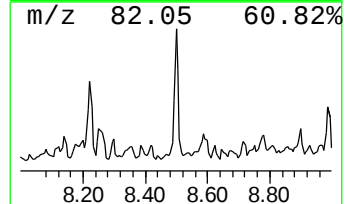
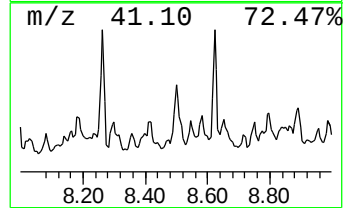
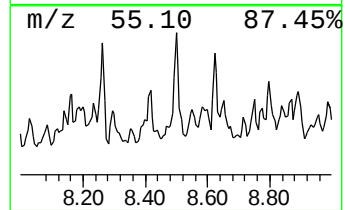
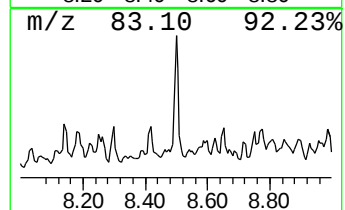
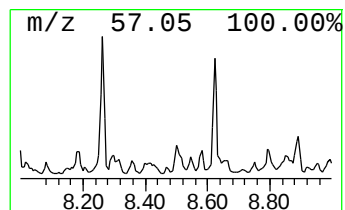
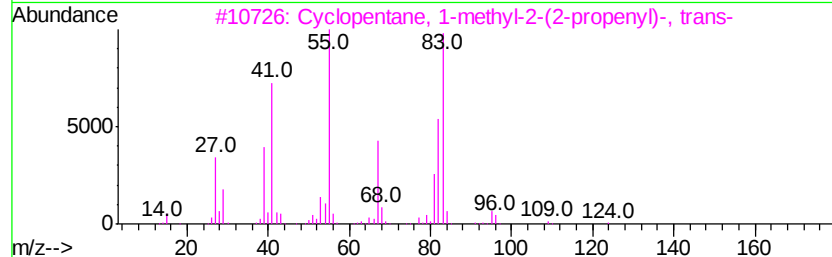
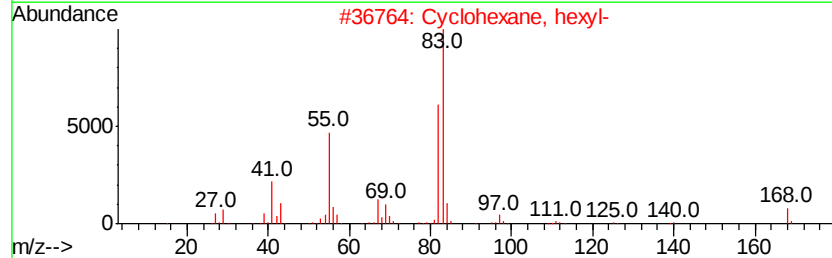
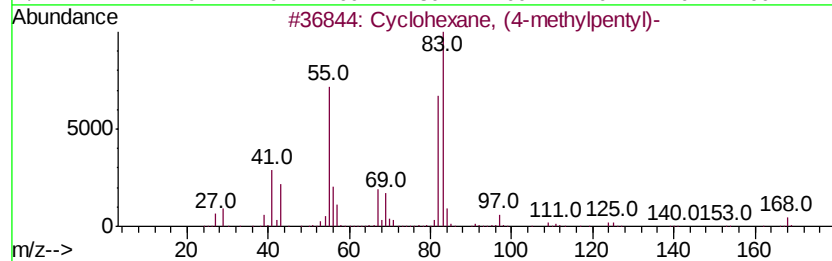
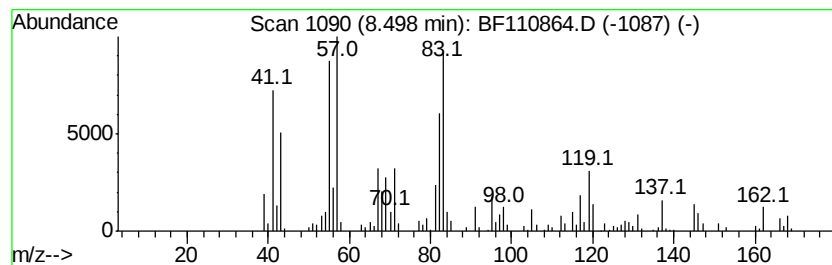
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown8.50 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.50	17.56 ng	347333	Naphthalene-d8	8.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	49
2		Cyclohexane, hexyl-	168	C12H24	004292-75-5	46
3		Cyclopentane, 1-methyl-2-(2-propenyl)-	124	C9H16	050746-53-7	46
4		Cyclohexane, propyl-	126	C9H18	001678-92-8	43
5		Cyclohexane, octyl-	196	C14H28	001795-15-9	43



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Acq On : 19 Nov 2018 20:15
Operator : JU/SJ
Sample : J5993-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

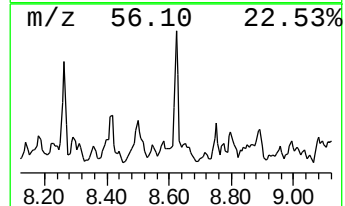
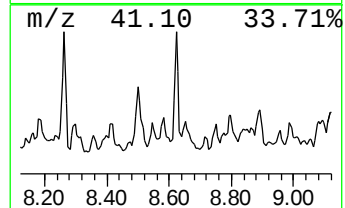
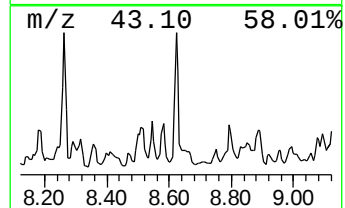
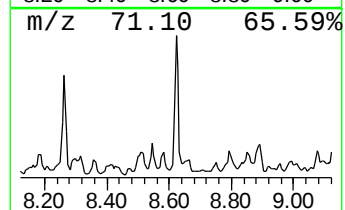
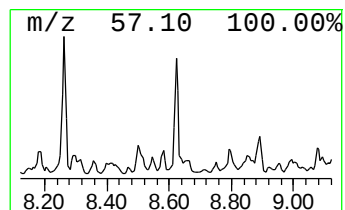
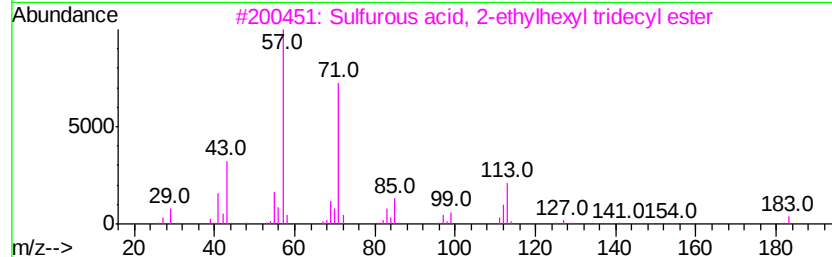
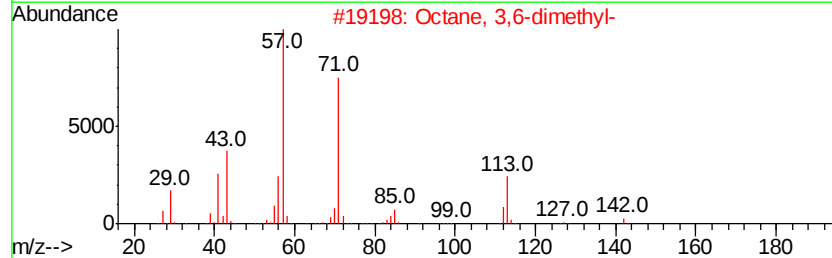
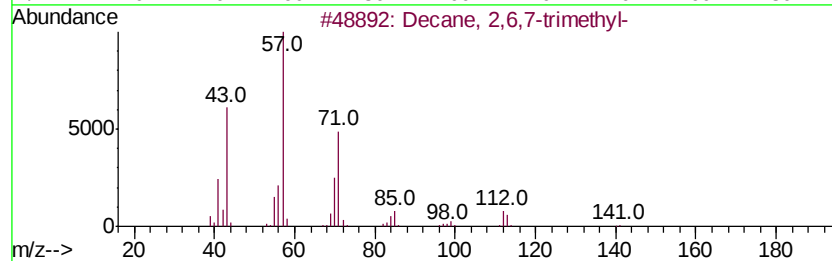
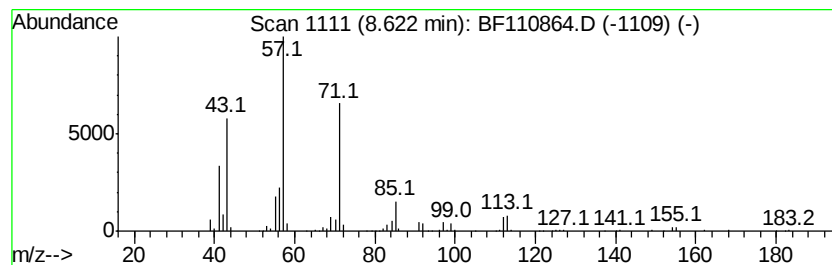
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Decane, 2,6,7-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
8.62	13.36 ng	264311	Naphthalene-d8	8.22		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 2,6,7-trimethyl-		184	C13H28	062108-25-2	78
2	Octane, 3,6-dimethyl-		142	C10H22	015869-94-0	64
3	Sulfurous acid, 2-ethylhexyl tri...		376	C21H44O3S	1000309-19-6	64
4	Dodecane, 2,6,11-trimethyl-		212	C15H32	031295-56-4	59
5	Pentadecane, 7-methyl-		226	C16H34	006165-40-8	53



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Data File : BF110864.D
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Misc :
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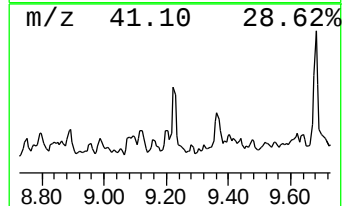
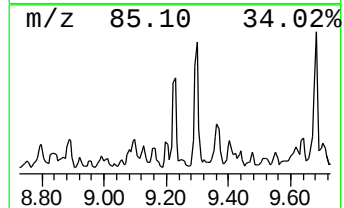
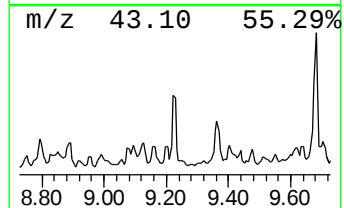
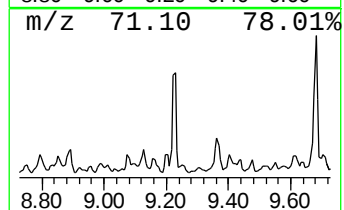
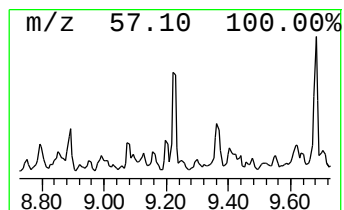
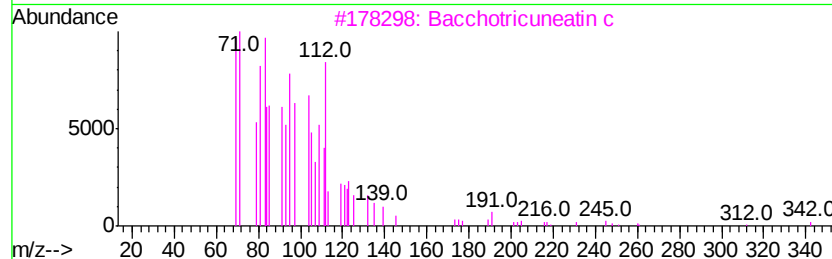
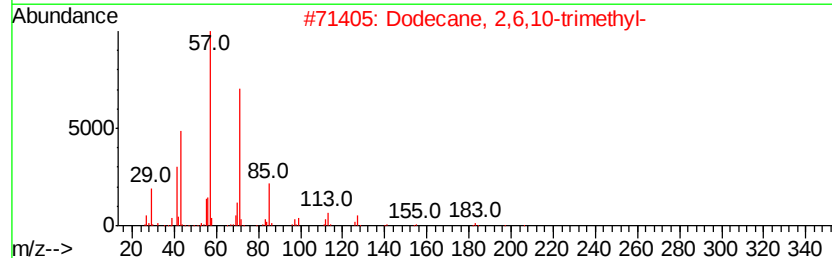
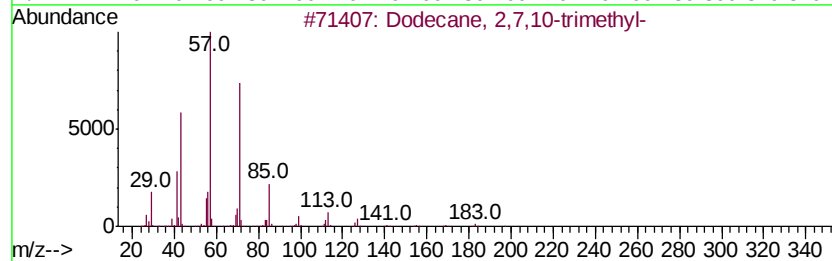
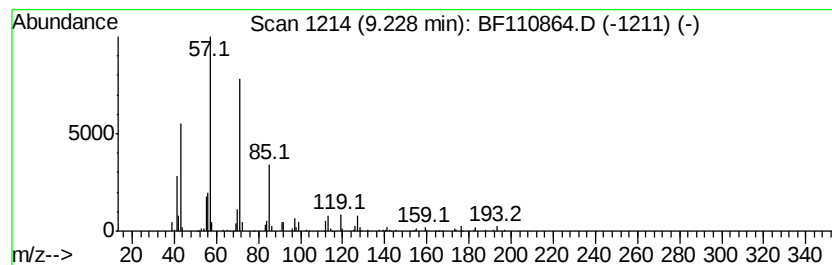
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Dodecane, 2,7,10-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.23	13.16 ng	239837	Acenaphthene-d10	9.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	90
2		Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	83
3		Bacchotricuneatin c	342	C20H22O5	066563-30-2	81
4		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72
5		Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	64



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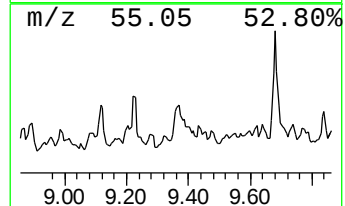
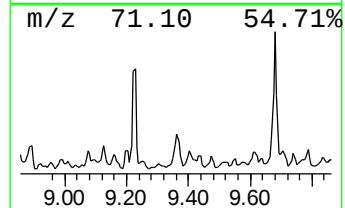
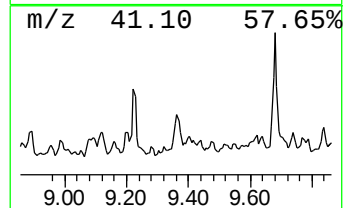
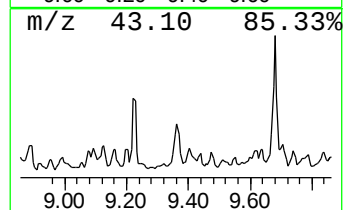
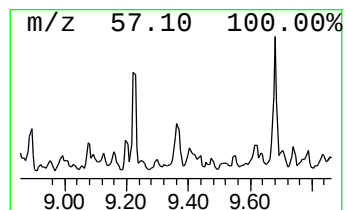
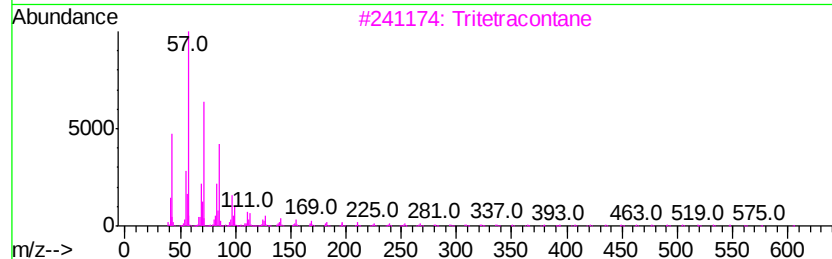
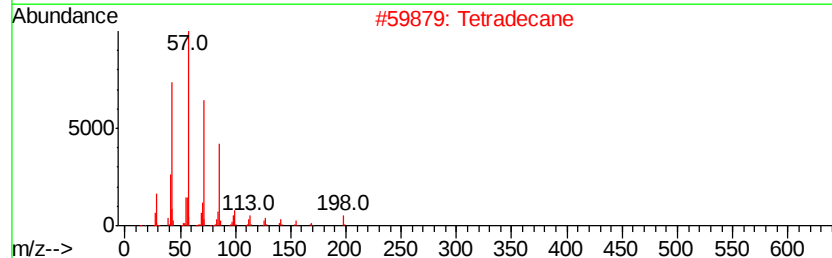
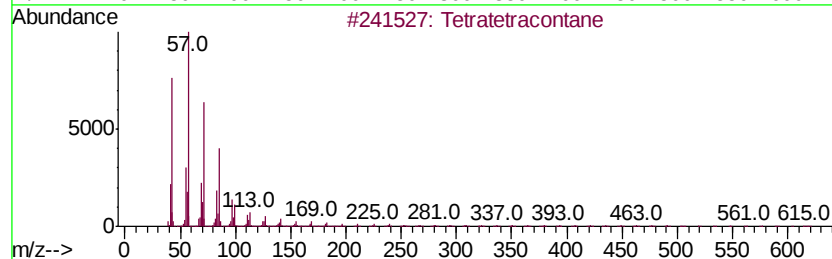
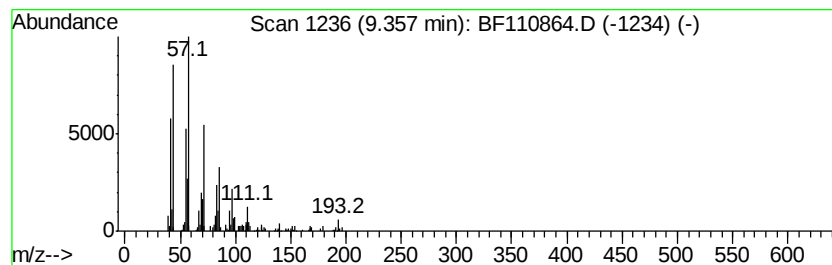
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Tetratetracontane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.36	14.64 ng	266803	Acenaphthene-d10	9.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	64
2		Tetradecane	198	C14H30	000629-59-4	55
3		Tritetracontane	605	C43H88	007098-21-7	52
4		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	52
5		Decane, 2,3,5,8-tetramethyl-	198	C14H30	192823-15-7	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
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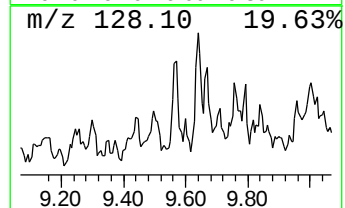
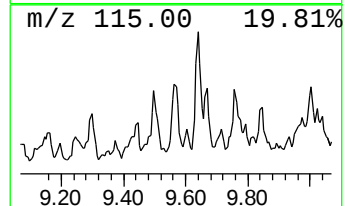
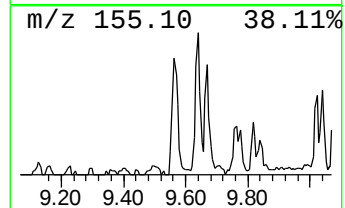
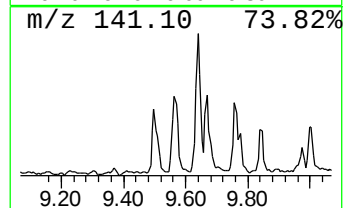
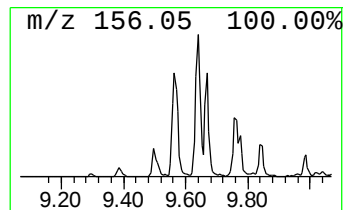
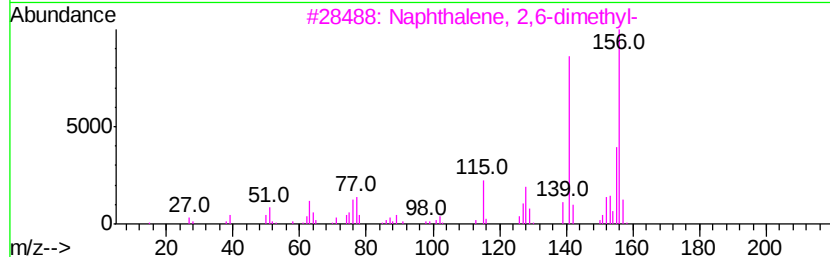
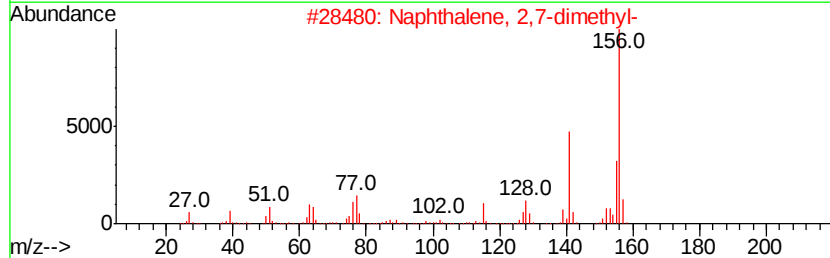
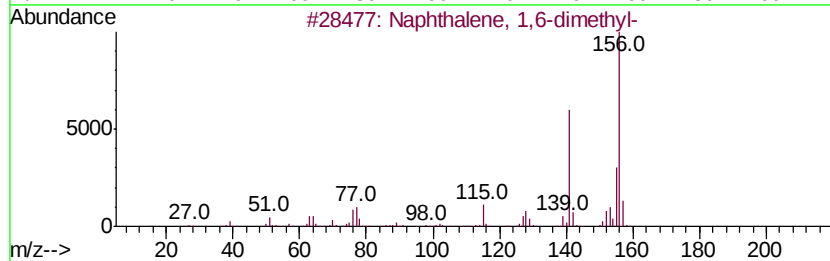
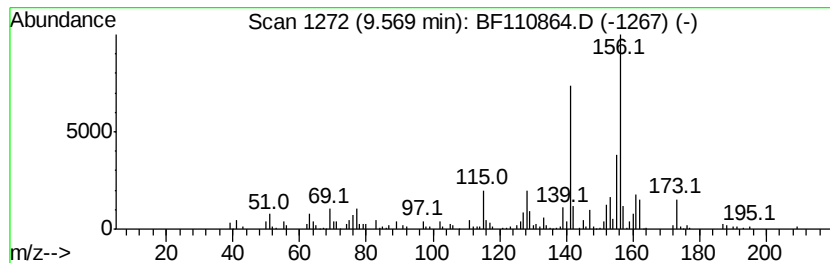
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1,6-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.57	11.90 ng	216843	Acenaphthene-d10	9.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
2		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
3		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	90
5		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	90



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Misc :
ALS Vial : 24 Sample Multiplier: 1

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ClientSampleId :
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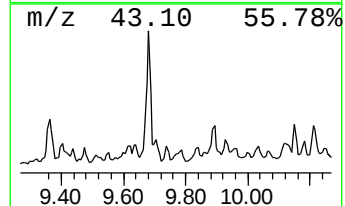
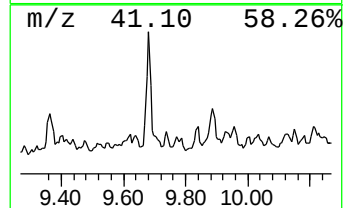
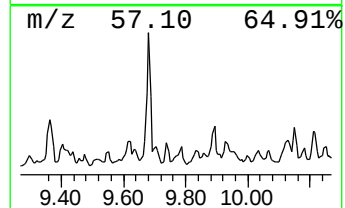
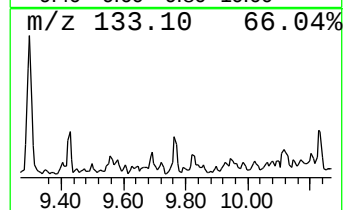
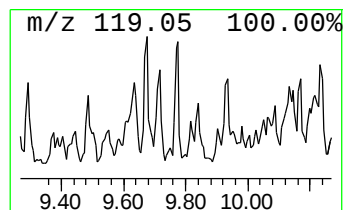
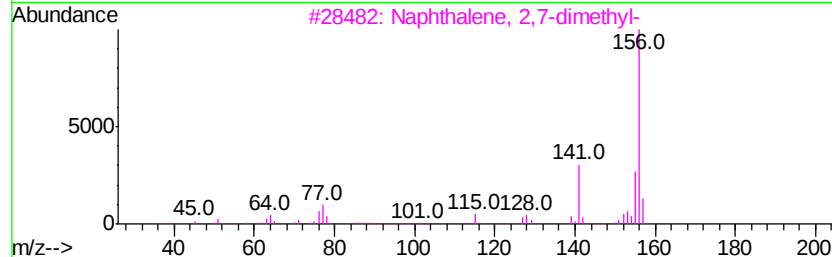
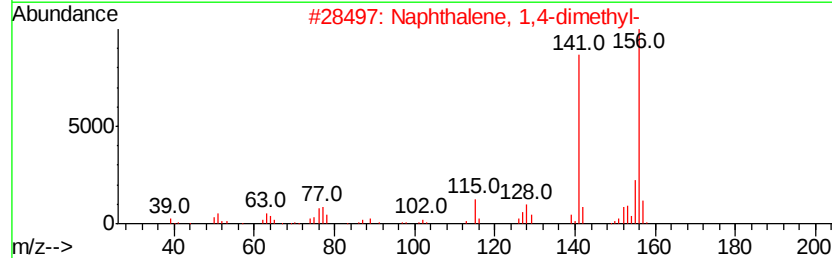
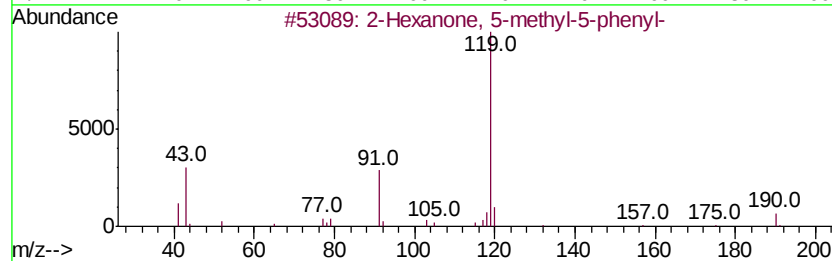
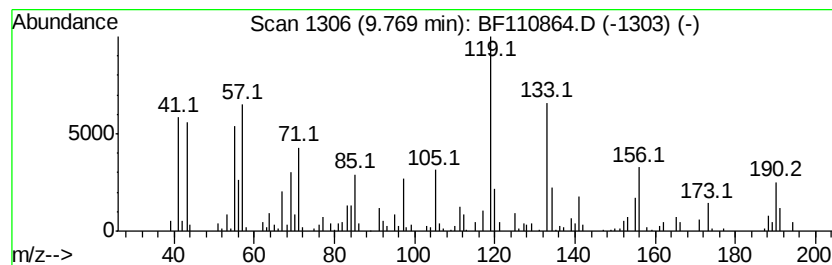
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown9.77 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.77	11.95 ng	217686	Acenaphthene-d10	9.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Hexanone, 5-methyl-5-phenyl-	190	C13H18O	014128-61-1	22
2		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	18
3		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	18
4		Propane, 2-cyclohexyl-2-phenyl-	202	C15H22	025683-97-0	14
5		Benzene, 1,4-dimethyl-2-(2-methy...	162	C12H18	055669-88-0	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110864.D
Acq On : 19 Nov 2018 20:15
Operator : JU/SJ
Sample : J5993-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20180929-AMENDED-WCC-2-COMP

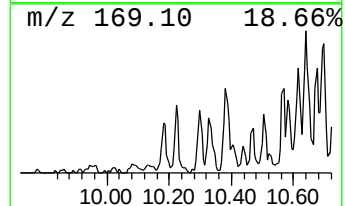
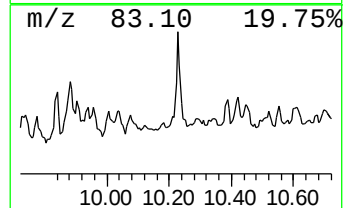
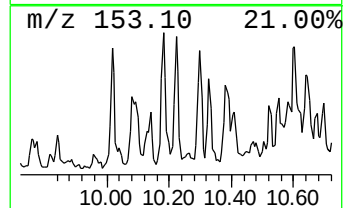
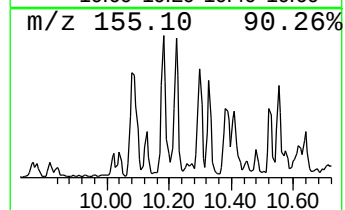
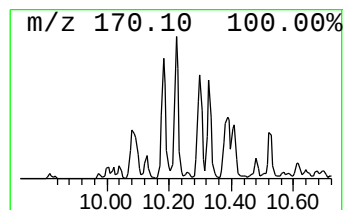
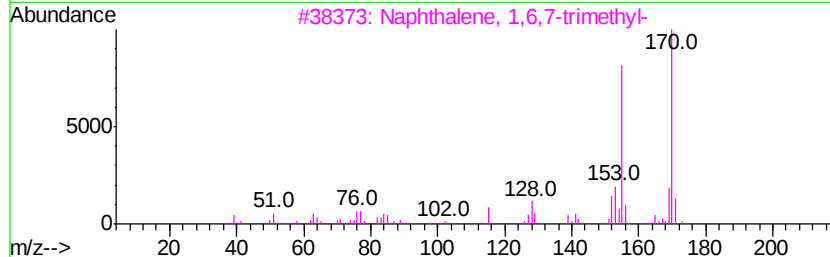
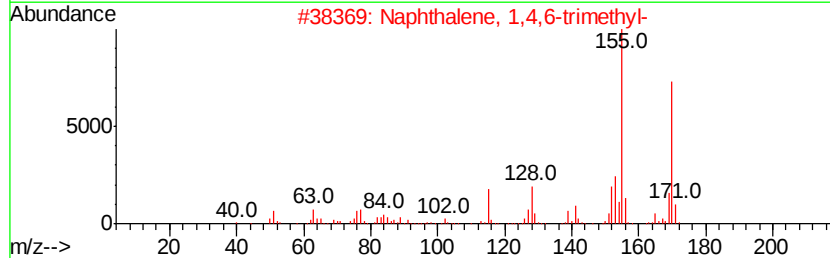
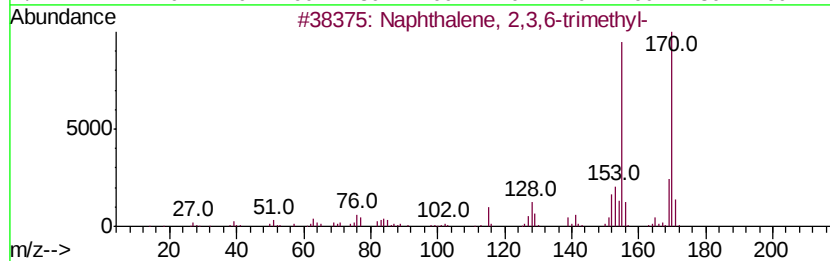
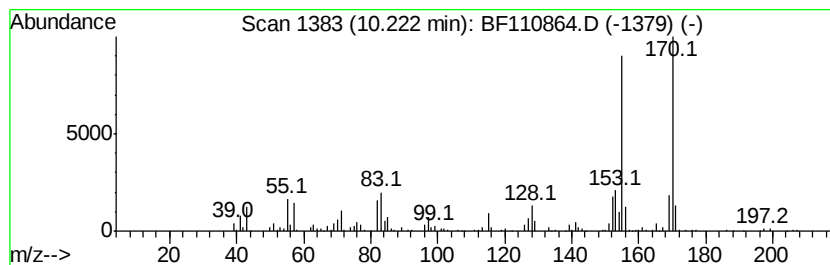
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Naphthalene, 2,3,6-trimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.22	12.12 ng	220770	Acenaphthene-d10	9.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	98
2		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
3		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
4		Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	91
5		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110864.D
Acq On : 19 Nov 2018 20:15
Operator : JU/SJ
Sample : J5993-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20180929-AMENDED-WCC-2-COMP

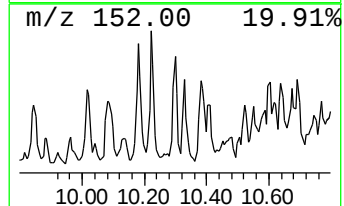
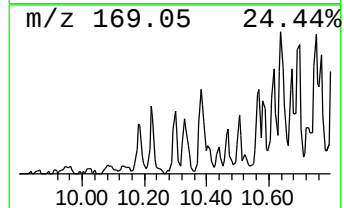
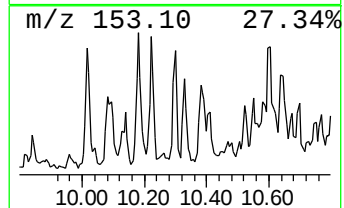
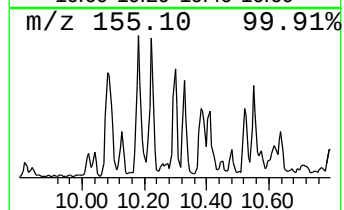
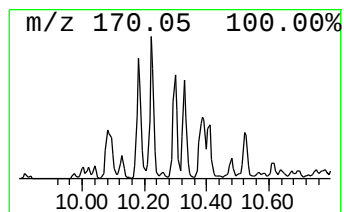
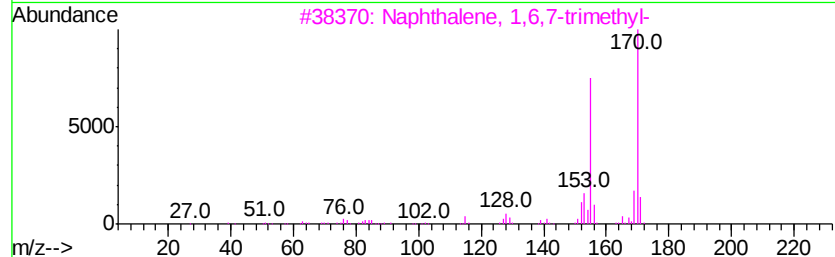
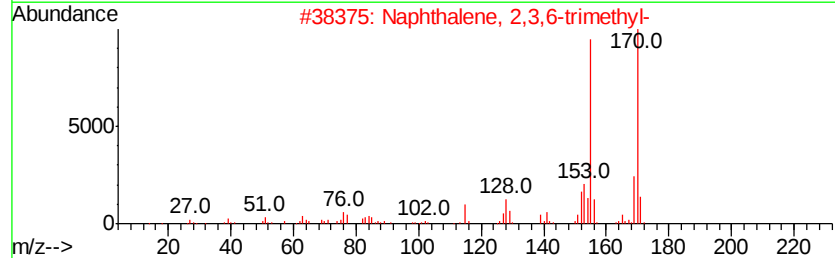
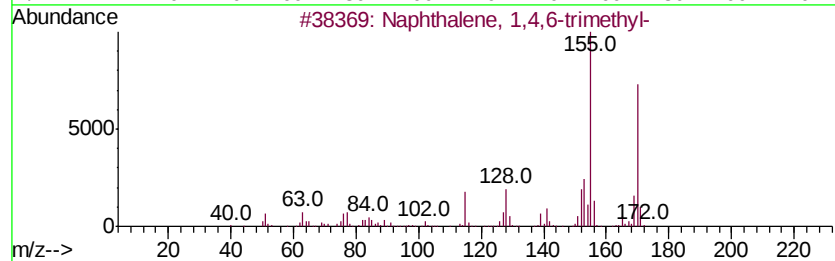
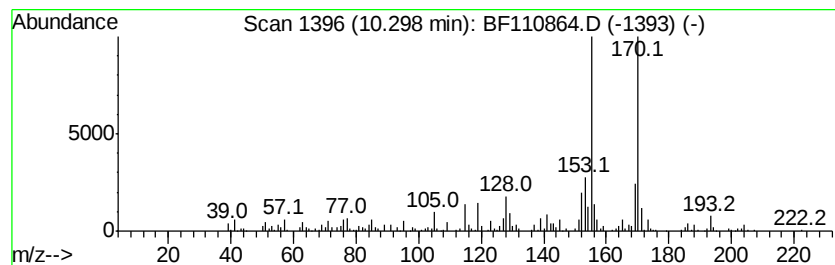
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Naphthalene, 1,4,6-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	12.41 ng	226064	Acenaphthene-d10	9.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	97
2		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
3		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	97
4		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	93
5		Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110864.D
Acq On : 19 Nov 2018 20:15
Operator : JU/SJ
Sample : J5993-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20180929-AMENDED-WCC-2-COMP

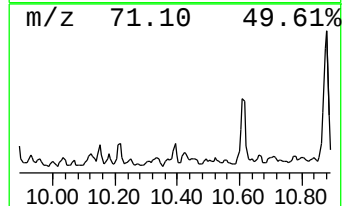
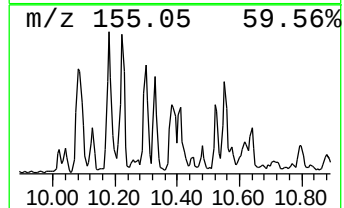
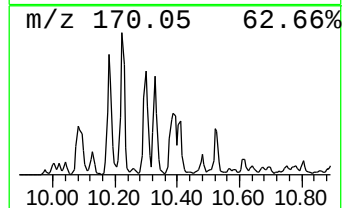
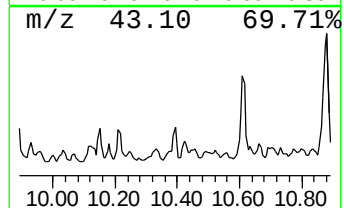
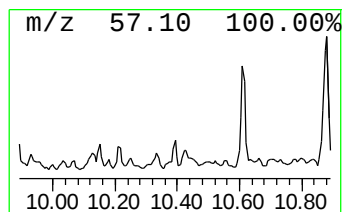
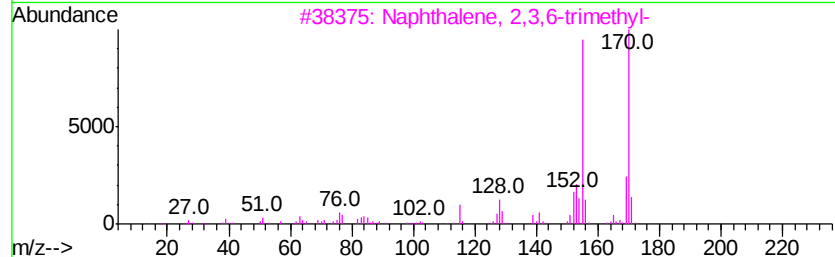
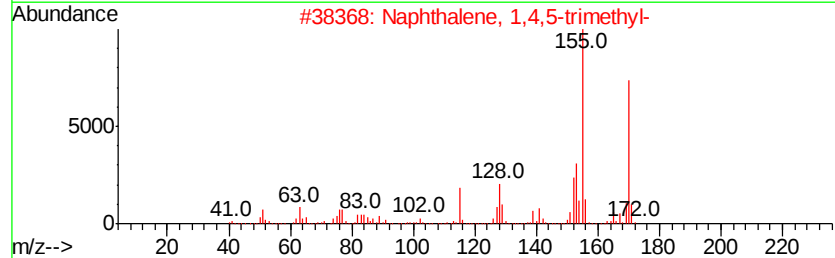
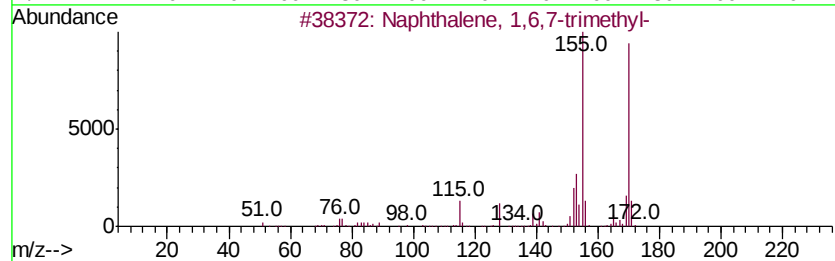
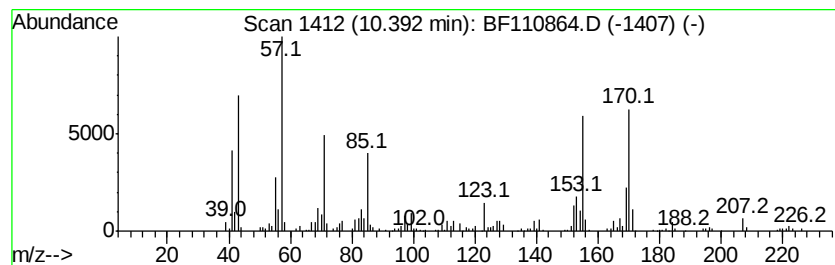
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Naphthalene, 1,6,7-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.39	15.24 ng	277628	Acenaphthene-d10	9.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	90
2		Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	86
3		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	78
4		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	55
5		1-(2,2-Dimethylcyclopropyl)-2-ph...	170	C13H14	1000294-91-1	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110864.D
Acq On : 19 Nov 2018 20:15
Operator : JU/SJ
Sample : J5993-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
20180929-AMENDED-WCC-2-COMP

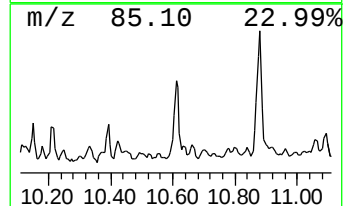
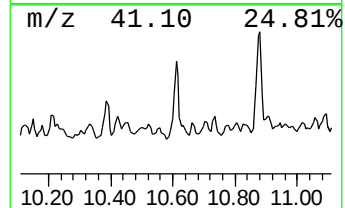
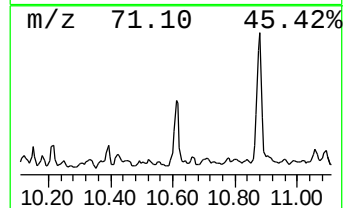
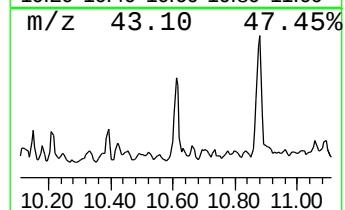
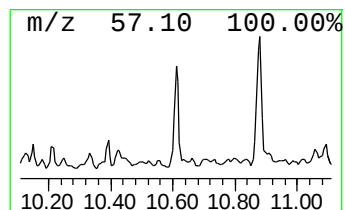
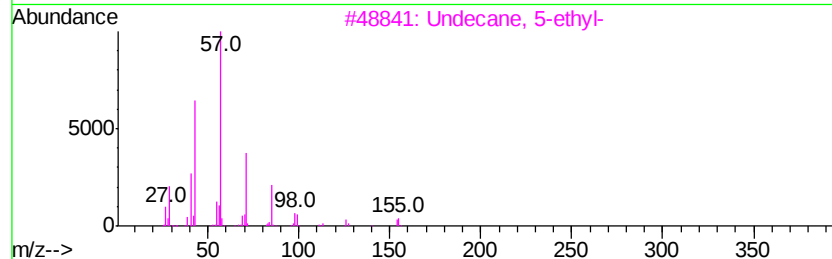
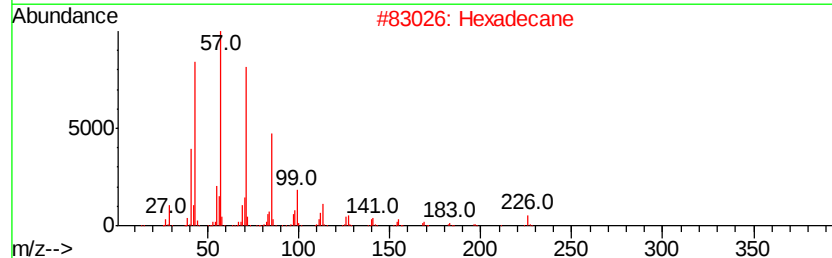
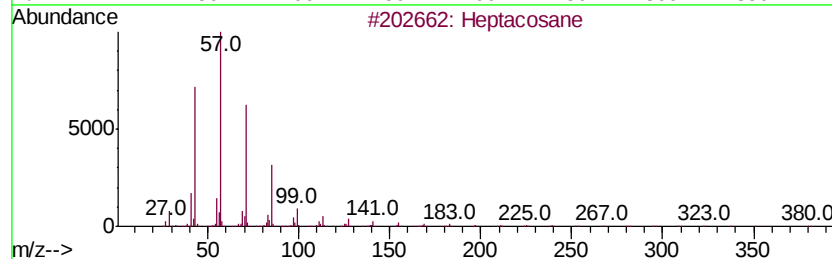
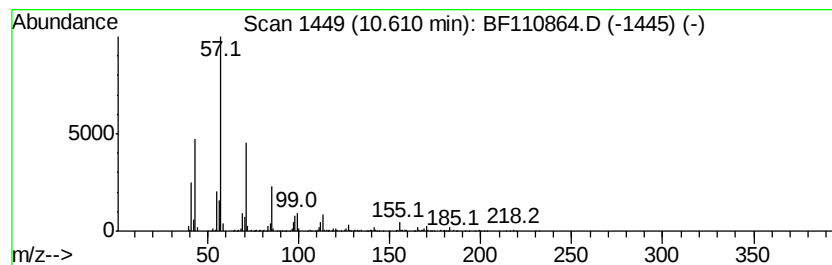
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Heptacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.61	26.46 ng	482105	Acenaphthene-d10	9.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	86
2		Hexadecane	226	C16H34	000544-76-3	80
3		Undecane, 5-ethyl-	184	C13H28	017453-94-0	78
4		Heneicosane	296	C21H44	000629-94-7	68
5		Eicosane	282	C20H42	000112-95-8	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110864.D
Acq On : 19 Nov 2018 20:15
Operator : JU/SJ
Sample : J5993-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

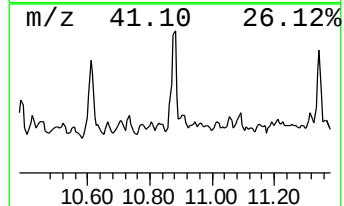
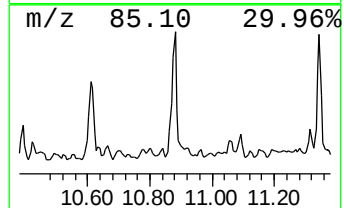
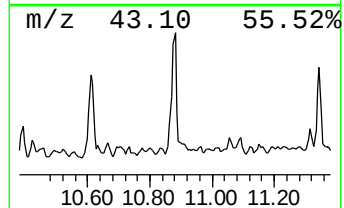
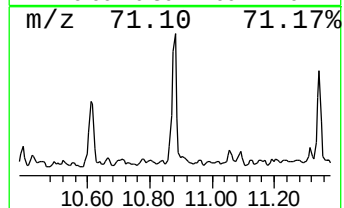
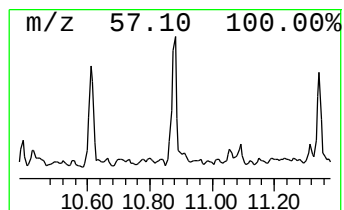
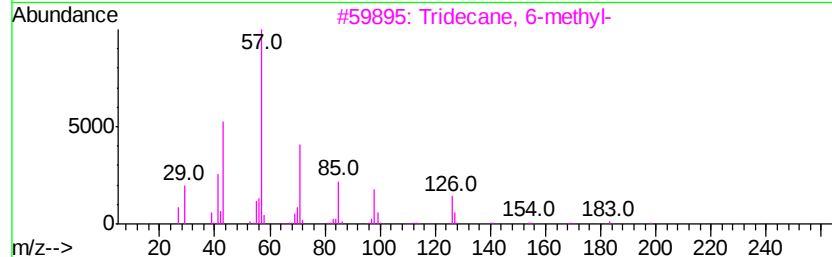
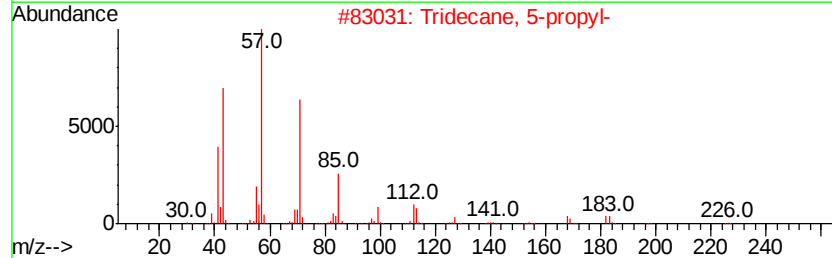
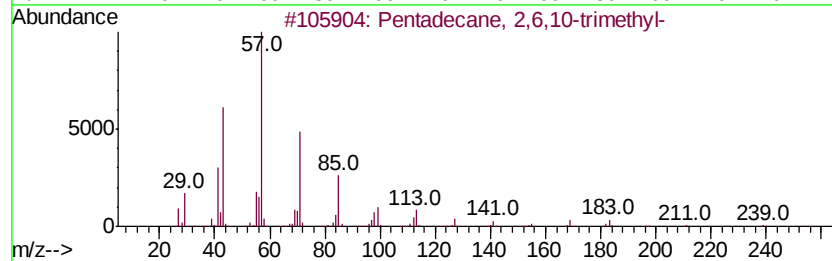
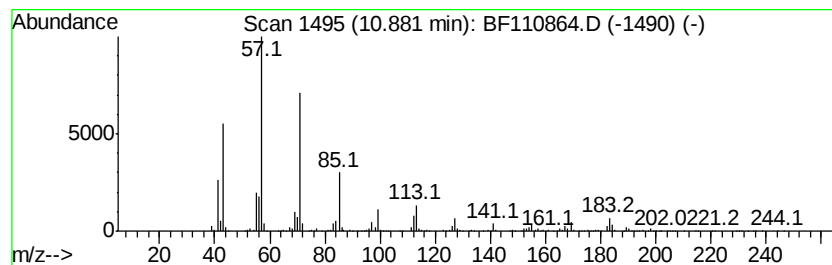
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ClientSampleId :
20180929-AMENDED-WCC-2-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Pentadecane, 2,6,10-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.88	56.73 ng	896899	Phenanthrene-d10		11.48
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2	Tridecane, 5-propyl-	226	C16H34	055045-11-9	86
3	Tridecane, 6-methyl-	198	C14H30	013287-21-3	81
4	Hexadecane, 2,6,10-trimethyl-	268	C19H40	055000-52-7	80
5	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110864.D
Acq On : 19 Nov 2018 20:15
Operator : JU/SJ
Sample : J5993-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
20180929-AMENDED-WCC-2-COMP

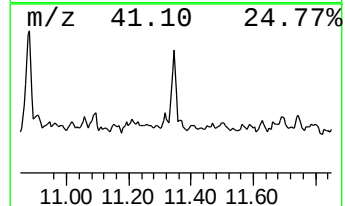
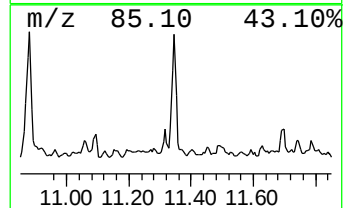
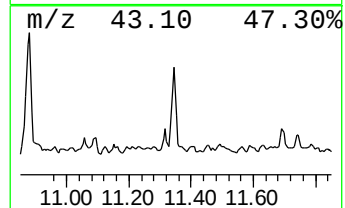
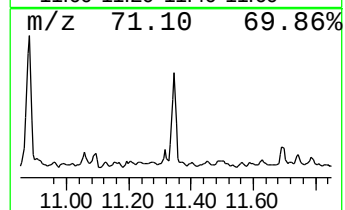
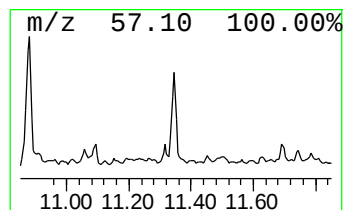
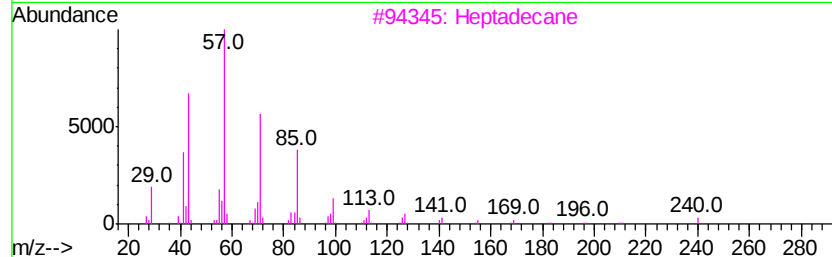
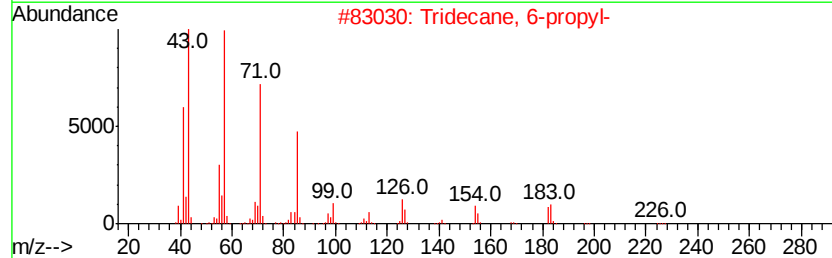
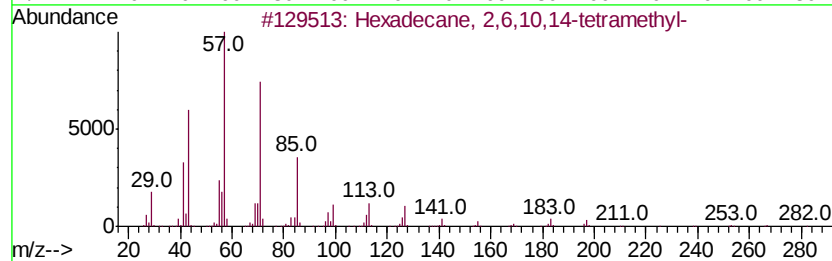
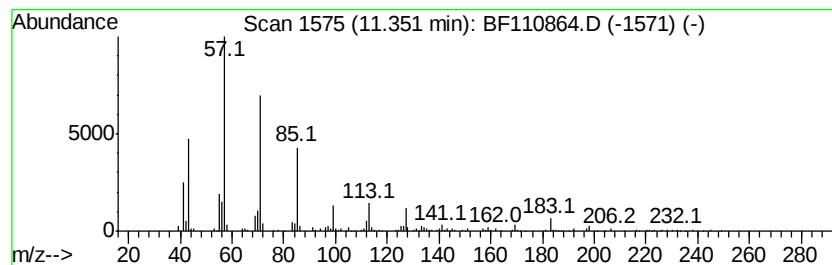
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Hexadecane, 2,6,10,14-tetra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.35	21.42 ng	338687	Phenanthrene-d10	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	91
2		Tridecane, 6-propyl-	226	C16H34	055045-10-8	90
3		Heptadecane	240	C17H36	000629-78-7	90
4		Heneicosane	296	C21H44	000629-94-7	86
5		Tridecane, 2-methyl-	198	C14H30	001560-96-9	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110864.D
Acq On : 19 Nov 2018 20:15
Operator : JU/SJ
Sample : J5993-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

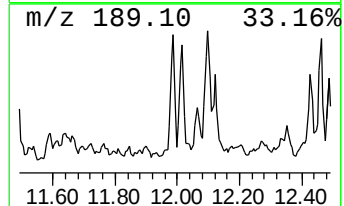
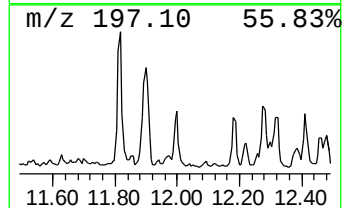
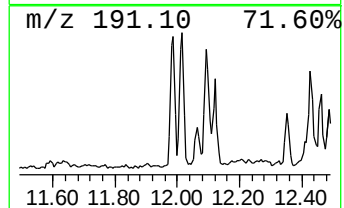
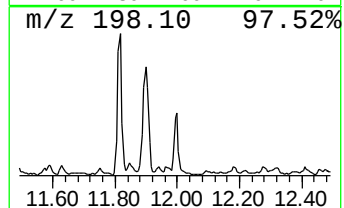
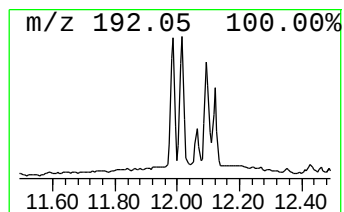
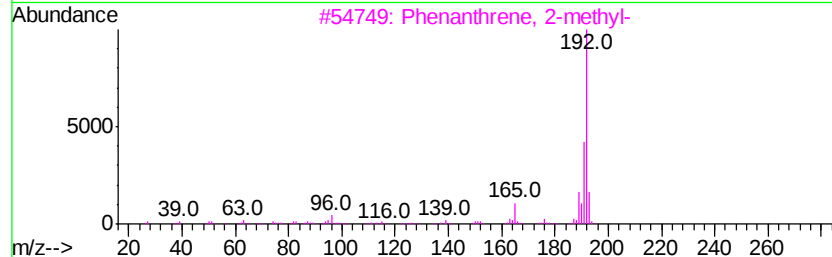
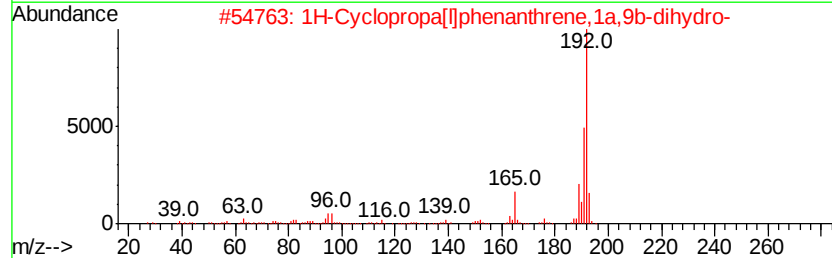
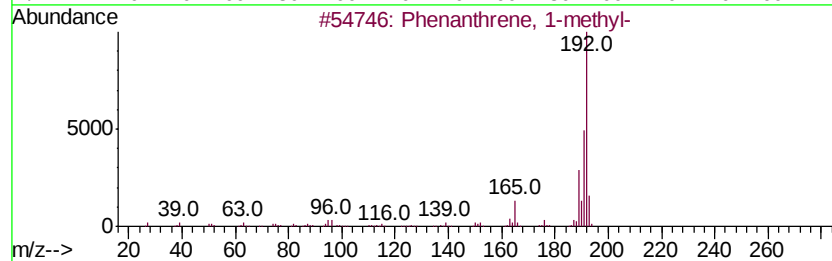
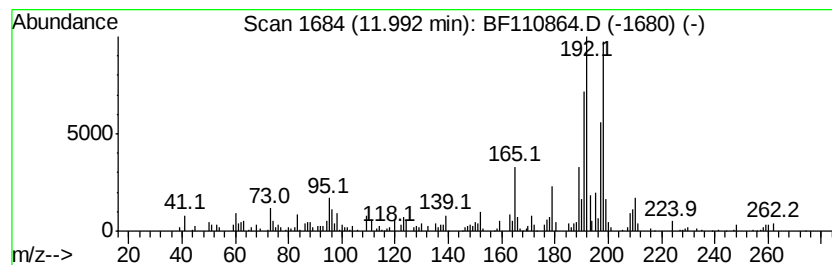
Instrument :
BNA_F
ClientSampleId :
20180929-AMENDED-WCC-2-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 Phenanthrene, 1-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.99	12.62 ng	199461	Phenanthrene-d10		11.48	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	96
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110864.D
Acq On : 19 Nov 2018 20:15
Operator : JU/SJ
Sample : J5993-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20180929-AMENDED-WCC-2-COMP

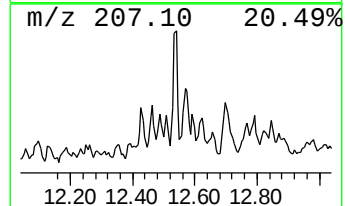
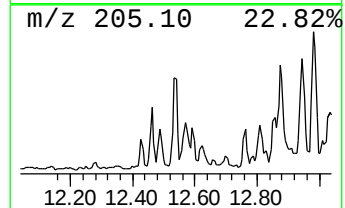
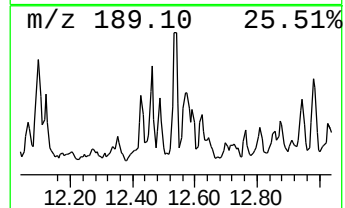
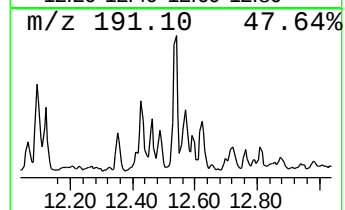
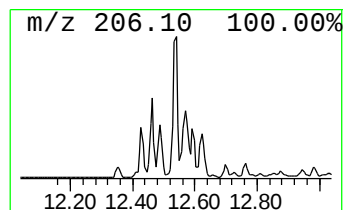
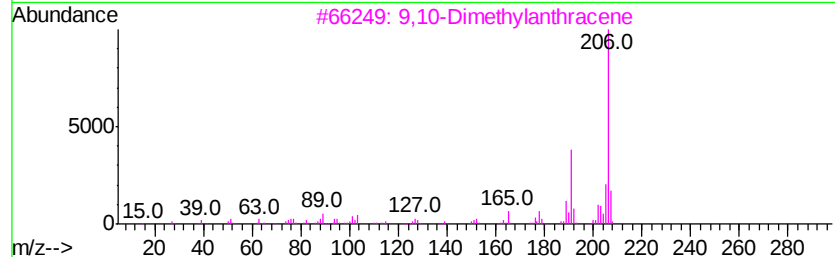
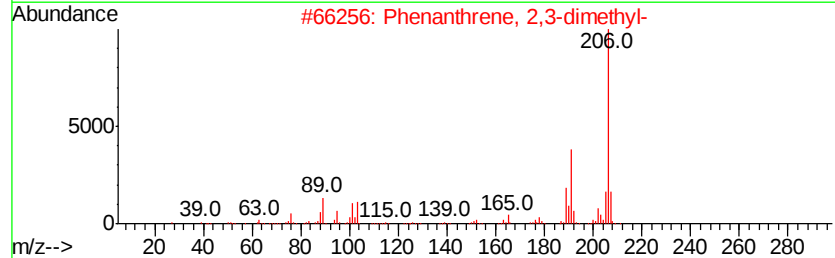
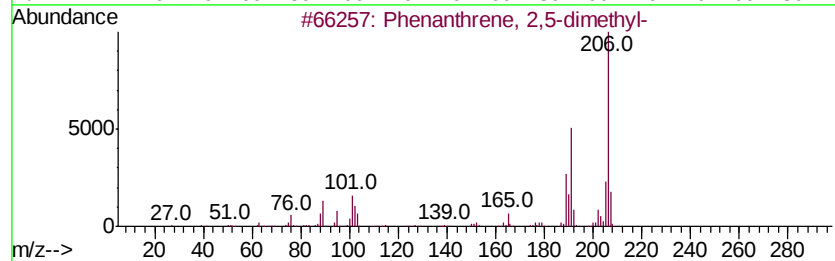
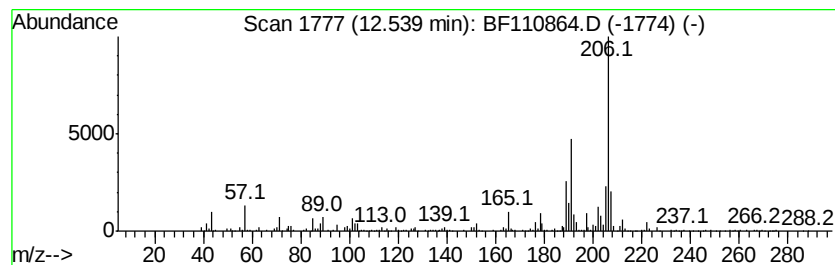
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Phenanthrene, 2,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	17.36 ng	274448	Phenanthrene-d10	11.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	94
2			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90
3			9,10-Dimethylanthracene	206	C16H14	000781-43-1	90
4			di-p-Tolylacetylene	206	C16H14	002789-88-0	90
5			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110864.D
Acq On : 19 Nov 2018 20:15
Operator : JU/SJ
Sample : J5993-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20180929-AMENDED-WCC-2-COMP

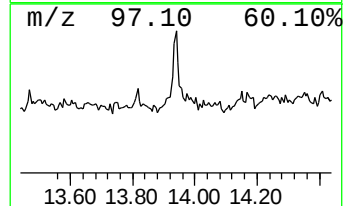
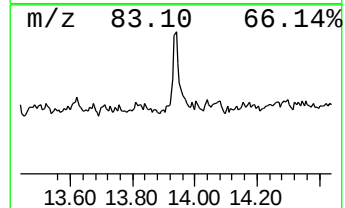
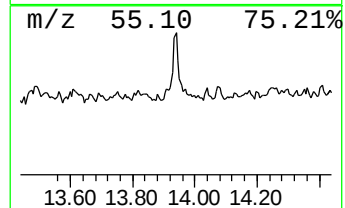
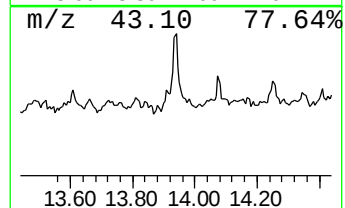
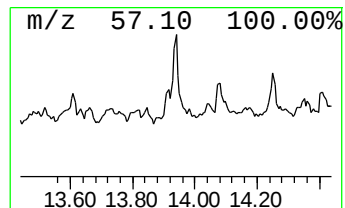
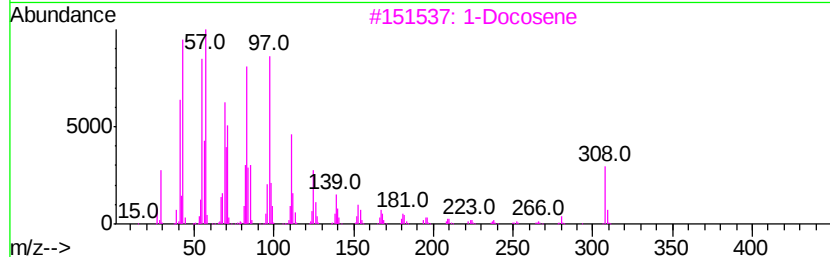
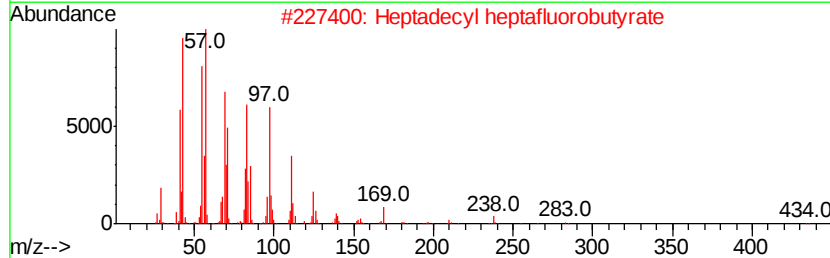
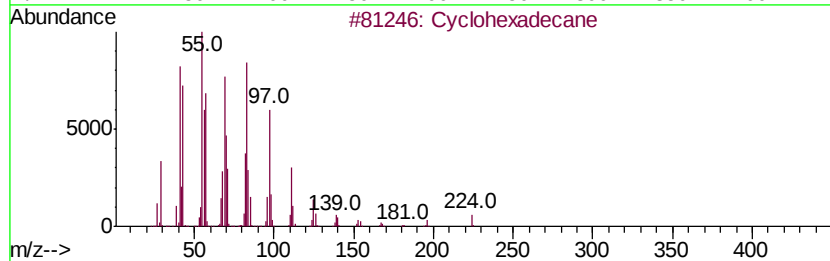
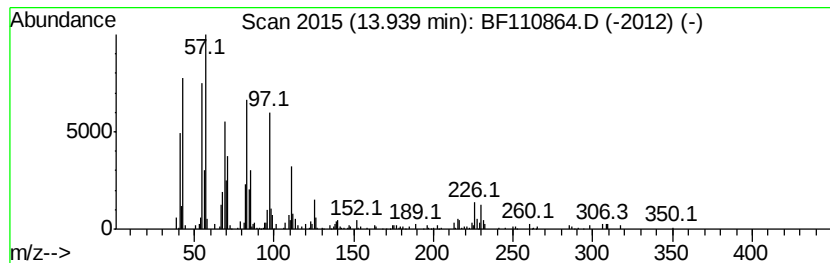
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 Cyclohexadecane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.94	11.40 ng	197493	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexadecane	224	C16H32	000295-65-8	95
2		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
3		1-Docosene	308	C22H44	001599-67-3	89
4		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	87
5		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111918\
 Data File : BF110864.D
 Acq On : 19 Nov 2018 20:15
 Operator : JU/SJ
 Sample : J5993-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20180929-AMENDED-WCC-2-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	45.3	ng	1034830	1	6.94	457062	20.0
Octane, 2,6-dimet...	6.17	13.3	ng	304442	1	6.94	457062	20.0
unknown6.71	6.71	40.2	ng	919580	1	6.94	457062	20.0
n-Hexanesulphonyl...	7.20	11.7	ng	268085	1	6.94	457062	20.0
unknown8.50	8.50	17.6	ng	347333	2	8.22	395581	20.0
Decane, 2,6,7-tri...	8.62	13.4	ng	264311	2	8.22	395581	20.0
Dodecane, 2,7,10-...	9.23	13.2	ng	239837	3	9.99	364416	20.0
Tetratetracontane	9.36	14.6	ng	266803	3	9.99	364416	20.0
Naphthalene, 1,6-...	9.57	11.9	ng	216843	3	9.99	364416	20.0
unknown9.77	9.77	11.9	ng	217686	3	9.99	364416	20.0
Naphthalene, 2,3,...	10.22	12.1	ng	220770	3	9.99	364416	20.0
Naphthalene, 1,4,...	10.30	12.4	ng	226064	3	9.99	364416	20.0
Naphthalene, 1,6,...	10.39	15.2	ng	277628	3	9.99	364416	20.0
Heptacosane	10.61	26.5	ng	482105	3	9.99	364416	20.0
Pentadecane, 2,6,...	10.88	56.7	ng	896899	4	11.48	316214	20.0
Hexadecane, 2,6,1...	11.35	21.4	ng	338687	4	11.48	316214	20.0
Phenanthrene, 1-m...	11.99	12.6	ng	199461	4	11.48	316214	20.0
Phenanthrene, 2,5...	12.54	17.4	ng	274448	4	11.48	316214	20.0
Cyclohexadecane	13.94	11.4	ng	197493	5	14.14	346413	20.0